

EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	1118	(514/394).CCLS.	US-PGPUB; USPAT; EPO; DERWENT	OR	OFF	2006/08/17 11:53
L2	301	(548/310.1).CCLS.	US-PGPUB; USPAT; EPO; DERWENT	OR	OFF	2006/08/17 11:53
S1	2	("5935983").PN.	US-PGPUB; USPAT; EPO; DERWENT	OR	OFF	2006/08/17 10:16
S2	2	("6649616").PN.	US-PGPUB; USPAT; EPO; DERWENT	OR	OFF	2006/08/17 10:31
S3	2	("7074780").PN.	US-PGPUB; USPAT; EPO; DERWENT	OR	OFF	2006/08/17 11:53

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1600RXA

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 "Ask CAS" for self-help around the clock
NEWS 3 FEB 27 New STN AnaVist pricing effective March 1, 2006
NEWS 4 APR 04 STN AnaVist \$500 visualization usage credit offered
NEWS 5 MAY 10 CA/CAPLUS enhanced with 1900-1906 U.S. patent records
NEWS 6 MAY 11 KOREAPAT updates resume
NEWS 7 MAY 19 Derwent World Patents Index to be reloaded and enhanced
NEWS 8 MAY 30 IPC 8 Rolled-up Core codes added to CA/CAPLUS and
USPATFULL/USPAT2
NEWS 9 MAY 30 The F-Term thesaurus is now available in CA/CAPLUS
NEWS 10 JUN 02 The first reclassification of IPC codes now complete in
INPADOC
NEWS 11 JUN 26 TULSA/TULSA2 reloaded and enhanced with new search and
and display fields
NEWS 12 JUN 28 Price changes in full-text patent databases EPFULL and PCTFULL
NEWS 13 JUL 11 CHEMSAFE reloaded and enhanced
NEWS 14 JUL 14 FSTA enhanced with Japanese patents
NEWS 15 JUL 19 Coverage of Research Disclosure reinstated in DWPI
NEWS 16 AUG 09 INSPEC enhanced with 1898-1968 archive

NEWS EXPRESS JUNE 30 CURRENT WINDOWS VERSION IS V8.01b, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 26 JUNE 2006.

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS LOGIN Welcome Banner and News Items
NEWS IPC8 For general information regarding STN implementation of IPC 8
NEWS X25 X.25 communication option no longer available

Enter NEWS followed by the item number or name to see news on that
specific topic.

All use of STN is subject to the provisions of the STN Customer
agreement. Please note that this agreement limits use to scientific
research. Use for software development or design or implementation
of commercial gateways or other similar uses is prohibited and may
result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 10:01:30 ON 17 AUG 2006

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 10:01:35 ON 17 AUG 2006

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2006 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 16 AUG 2006 HIGHEST RN 902024-59-3
DICTIONARY FILE UPDATES: 16 AUG 2006 HIGHEST RN 902024-59-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

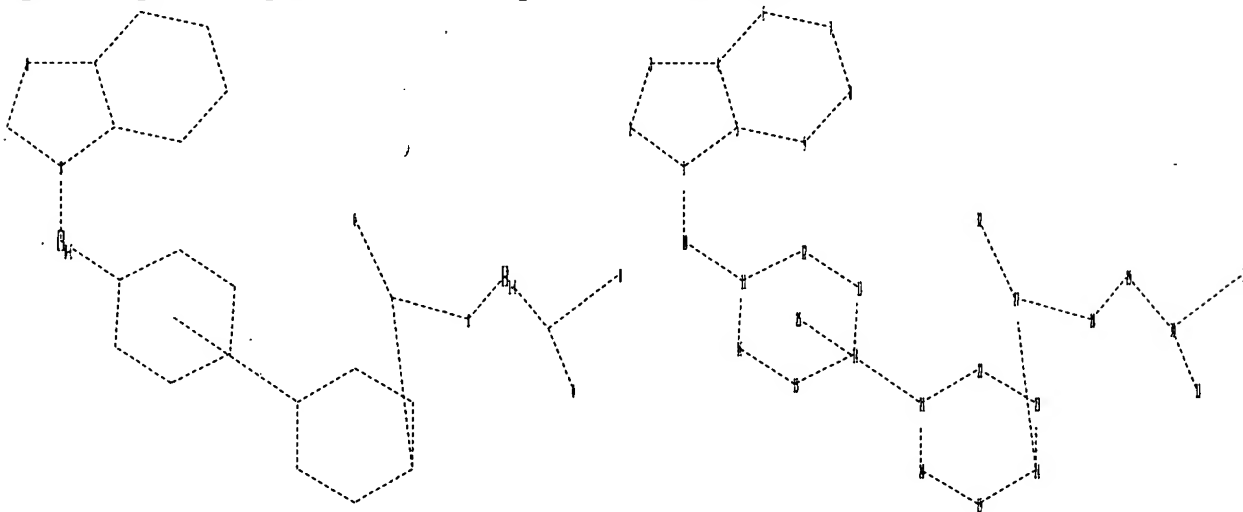
Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>

Uploading C:\Program Files\Stnexp\Queries\QUERIES\10805548.str



chain nodes :

10 27 28 29 30 31 32 33

ring nodes :

1 2 3 4 5 6 7 8 9 11 12 13 14 15 16 19 20 21 22 23 24

chain bonds :

1-10 10-11 24-27 27-32 27-28 28-29 29-30 30-31 30-33

ring bonds :

1-2 1-5 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 11-12 11-16 12-13 13-14 14-15
15-16 19-20 19-24 20-21 21-22 22-23 23-24

exact/norm bonds :

1-2 1-5 1-10 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 10-11 11-12 11-16 12-13
13-14 14-15 15-16 19-20 19-24 20-21 21-22 22-23 23-24 24-27 27-32 27-28
28-29 29-30 30-31 30-33

isolated ring systems :

containing 1 : 11 : 19 :

Match level :

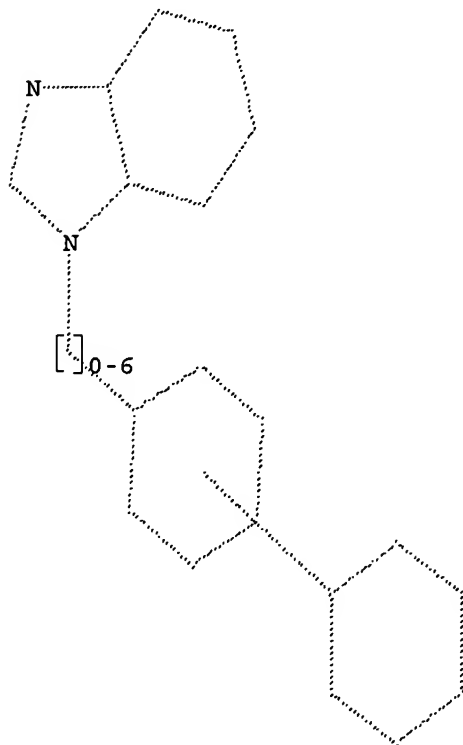
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 19:Atom 20:Atom 21:Atom
22:Atom 23:Atom 24:Atom 25:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS
31:CLASS 32:CLASS 33:CLASS

L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 10:02:48 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 5819 TO ITERATE

34.4% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

50 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**

PROJECTED ITERATIONS: 111807 TO 120953
PROJECTED ANSWERS: 2853 TO 4477

L2 50 SEA,SSS SAM L1

=> fil stnguide

COST IN U.S. DOLLARS

SINCE FILE
ENTRY
0.88

TOTAL
SESSION
1.09

FULL ESTIMATED COST

FILE 'STNGUIDE' ENTERED AT 10:02:55 ON 17 AUG 2006
USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT
COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY, JAPAN SCIENCE

AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.

LAST RELOADED: Aug 11, 2006 (20060811/UP).

=>

Uploading

THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE

Do you want to switch to the Registry File?

Choice (Y/n):

Switching to the Registry File...

Some commands only work in certain files. For example, the EXPAND command can only be used to look at the index in a file which has an index. Enter "HELP COMMANDS" at an arrow prompt (=>) for a list of commands which can be used in this file.

=> FILE REGISTRY

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.12	1.21

FILE 'REGISTRY' ENTERED AT 10:04:23 ON 17 AUG 2006

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2006 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 16 AUG 2006 HIGHEST RN 902024-59-3

DICTIONARY FILE UPDATES: 16 AUG 2006 HIGHEST RN 902024-59-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 6, 2006

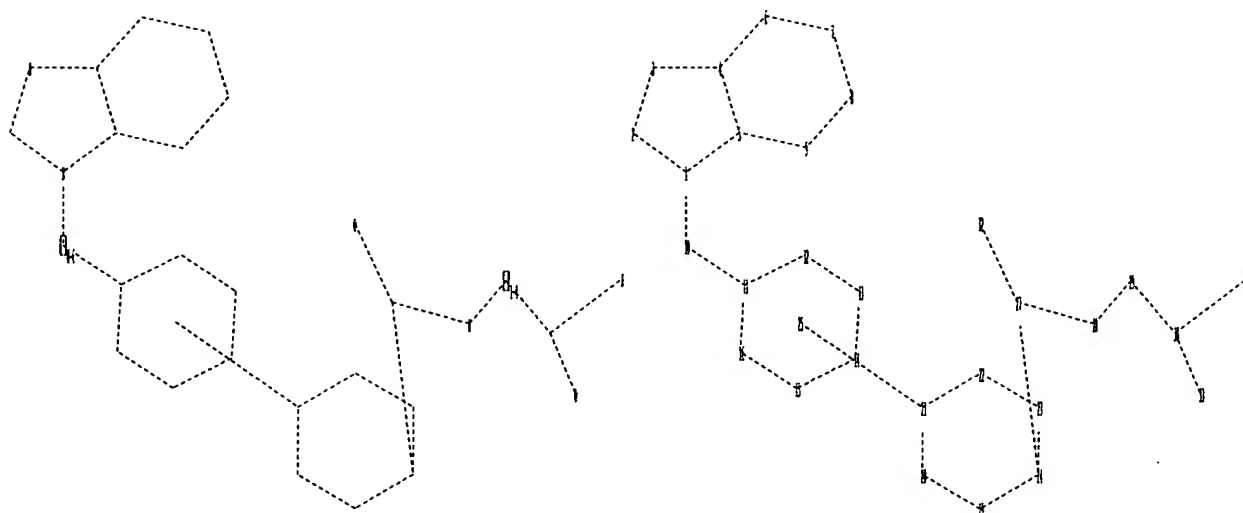
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>

Uploading C:\Program Files\Stnexp\Queries\QUERIES\10805548.str



```

chain nodes :
10 27 28 29 30 31 32 33
ring nodes :
1 2 3 4 5 6 7 8 9 11 12 13 14 15 16 19 20 21 22 23 24
chain bonds :
1-10 10-11 24-27 27-32 27-28 28-29 29-30 30-31 30-33
ring bonds :
1-2 1-5 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 11-12 11-16 12-13 13-14 14-15
15-16 19-20 19-24 20-21 21-22 22-23 23-24
exact/norm bonds :
1-2 1-5 1-10 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 10-11 11-12 11-16 12-13
13-14 14-15 15-16 19-20 19-24 20-21 21-22 22-23 23-24 24-27 27-32 27-28
28-29 29-30 30-31 30-33
isolated ring systems :
containing 1 : 11 : 19 :

```

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 19:Atom 20:Atom 21:Atom
22:Atom 23:Atom 24:Atom 25:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS
31:CLASS 32:CLASS 33:CLASS

```

L3 STRUCTURE UPLOADED

=> d

L3 HAS NO ANSWERS

L3 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> d his

(FILE 'HOME' ENTERED AT 10:01:30 ON 17 AUG 2006)

FILE 'REGISTRY' ENTERED AT 10:01:35 ON 17 AUG 2006

L1 STRUCTURE UPLOADED

L2 50 S L1

FILE 'STNGUIDE' ENTERED AT 10:02:55 ON 17 AUG 2006

FILE 'REGISTRY' ENTERED AT 10:04:23 ON 17 AUG 2006

L3 STRUCTURE UPLOADED

=> s l3

SAMPLE SEARCH INITIATED 10:04:44 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 67 TO ITERATE

100.0% PROCESSED 67 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 849 TO 1831

PROJECTED ANSWERS: 0 TO 0

L4 0 SEA SSS SAM L3

=> s l3 full

FULL SEARCH INITIATED 10:04:49 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 1388 TO ITERATE

100.0% PROCESSED 1388 ITERATIONS

23 ANSWERS

SEARCH TIME: 00.00.01

L5 23 SEA SSS FUL L3

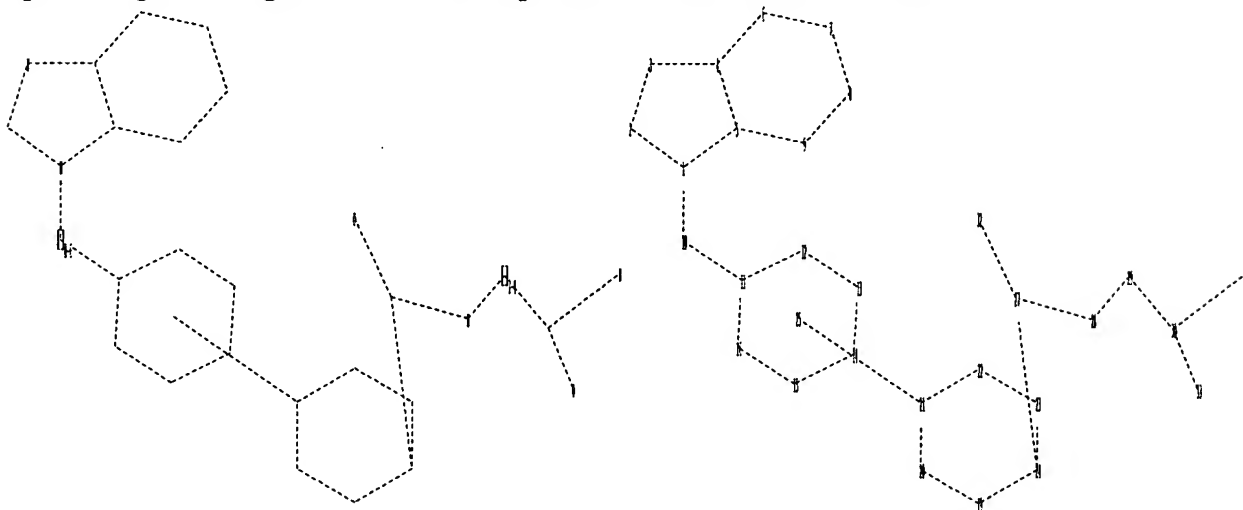
=> s l5 and caplus/lc

51672654 CAPLUS/LC

L6 23 L5 AND CAPLUS/LC

=>

Uploading C:\Program Files\Stnexp\Queries\QUERIES\10805548.str



chain nodes :

10 27 28 29 30 31 32 33

ring nodes :

1 2 3 4 5 6 7 8 9 11 12 13 14 15 16 19 20 21 22 23 24

chain bonds :

1-10 10-11 24-27 27-32 27-28 28-29 29-30 30-31 30-33

ring bonds :

1-2 1-5 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 11-12 11-16 12-13 13-14 14-15
15-16 19-20 19-24 20-21 21-22 22-23 23-24

exact/norm bonds :

1-2 1-5 1-10 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 10-11 11-12 11-16 12-13
13-14 14-15 15-16 19-20 19-24 20-21 21-22 22-23 23-24 24-27 27-32 27-28
28-29 29-30 30-31 30-33
isolated ring systems :
containing 1 : 11 : 19 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 19:Atom 20:Atom 21:Atom
22:Atom 23:Atom 24:Atom 25:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS
31:CLASS 32:CLASS 33:CLASS

L7 STRUCTURE UPLOADED

=> d

L7 HAS NO ANSWERS

L7 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s l7

SAMPLE SEARCH INITIATED 10:05:32 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 67 TO ITERATE

100.0% PROCESSED 67 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 849 TO 1831

PROJECTED ANSWERS: 0 TO 0

L8 0 SEA SSS SAM L7

=> s l7 full

FULL SEARCH INITIATED 10:05:35 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 1388 TO ITERATE

100.0% PROCESSED 1388 ITERATIONS

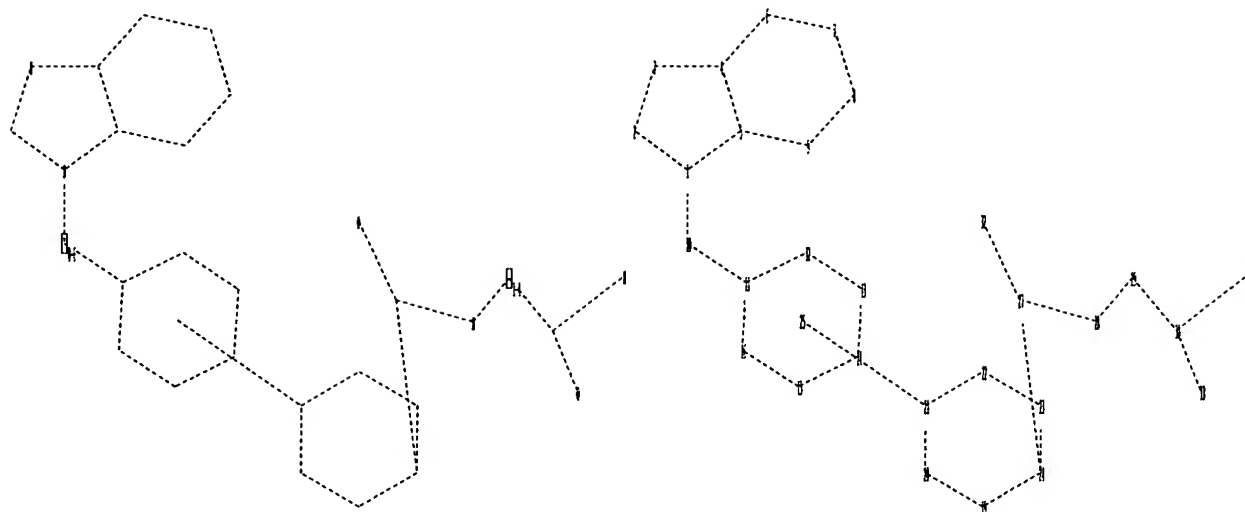
0 ANSWERS

SEARCH TIME: 00.00.01

L9 0 SEA SSS FUL L7

=>

Uploading C:\Program Files\Stnexp\Queries\QUERIES\10805548.str



```

chain nodes :
10 27 28 29 30 31 32 33
ring nodes :
1 2 3 4 5 6 7 8 9 11 12 13 14 15 16 19 20 21 22 23 24
chain bonds :
1-10 10-11 24-27 27-32 27-28 28-29 29-30 30-31 30-33
ring bonds :
1-2 1-5 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 11-12 11-16 12-13 13-14 14-15
15-16 19-20 19-24 20-21 21-22 22-23 23-24
exact/norm bonds :
1-2 1-5 1-10 2-3 3-4 4-5 4-6 5-9 6-7 7-8 8-9 10-11 11-12 11-16 12-13
13-14 14-15 15-16 19-20 19-24 20-21 21-22 22-23 23-24 24-27 27-32 27-28
28-29 29-30 30-31 30-33
isolated ring systems :
containing 1 : 11 : 19 :

```

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 19:Atom 20:Atom 21:Atom
22:Atom 23:Atom 24:Atom 25:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS
31:CLASS 32:CLASS 33:CLASS

```

L10 STRUCTURE UPLOADED

=> d

L10 HAS NO ANSWERS

L10 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s l10

SAMPLE SEARCH INITIATED 10:06:09 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 67 TO ITERATE

100.0% PROCESSED 67 ITERATIONS

SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 849 TO 1831
PROJECTED ANSWERS: 0 TO 0

L11 0 SEA SSS SAM L10

=> s l10 full

FULL SEARCH INITIATED 10:06:11 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1388 TO ITERATE

100.0% PROCESSED 1388 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

L12 0 SEA SSS FUL L10

=> fil caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	505.58	506.79

FILE 'CAPLUS' ENTERED AT 10:06:14 ON 17 AUG 2006
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 17 Aug 2006 VOL 145 ISS 8
FILE LAST UPDATED: 16 Aug 2006 (20060816/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/infopolicy.html>

=> s l6

L13 3 L6

=> d ibib abs hitstr 1-3

ACCESSION NUMBER: 2000:841902 CAPLUS
DOCUMENT NUMBER: 133:362969TITLE: Synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cardiovascular ischemia
INVENTOR(S): Bischoff, Erwin; Lensky, Stephan; Muller, Stephan
Nicholas; Paulsen, Holger; Keldenich, Jorg; Krahn, Thomas; Schuhmacher, JoachimPATENT ASSIGNEE(S): Bayer A.-G., Germany
SOURCE: Ger. Offen., 30 pp.

CODEN: GWXXBX

DOCUMENT TYPE: Patent

LANGUAGE: German

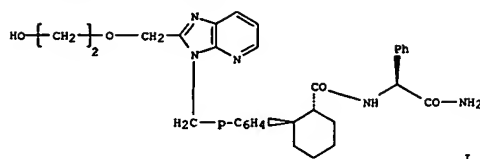
FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 19924819	A1	20001130	DE 1999-19924819	19990529
CA 2375188	AA	20001207	CA 2000-2375188	20000516
WO 2000073274	A2	20001207	WO 2000-EP4431	20000516
WO 2000073274	A3	20010419		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW				
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
BR 2000011049	A	20020319	BR 2000-11049	20000516
EP 1187812	A2	20020320	EP 2000-925290	20000516
EP 1187812	B1	20050907		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
TR 200103398	T2	20020321	TR 2001-3398	20000516
JP 2003500474	T2	20030107	JP 2000-621340	20000516
AU 766140	B2	20031009	AU 2000-44057	20000516
AT 303995	E	20050915	AT 2000-925290	20000516
ES 2248071	T3	20060316	ES 2000-925290	20000516
ZA 2001008894	A	20021029	ZA 2001-8894	20011029
US 6984642	B1	20060110	US 2001-980242	20011129
HK 1047107	A1	20050520	HK 2002-108818	20021204
			DE 1999-19924819	19990529
			WO 2000-EP4431	20000516

OTHER SOURCE(S): MARPAT 133:362969

GI



AB Title compds., e.g. (I), were prepared for use in treating cardiovascular ischemic disorders in humans or animals. Thus, 2-(2-hydroxyethoxymethyl)pyrido[2,3-d]imidazole (preparation given) was reacted with

(1R,2R)-2-(4-methylphenyl)cyclohexanecarboxylic acid (resolution from racemate given) to yield the intermediate material which was reacted with (S)-phenylglycineamide hydrochloride to give I. In in vitro tests of rabbit erythrocyte adenosine uptake, the 2-(morpholin-4-yl)methyl (in place of the 2-(2-hydroxyethoxymethyl) sidechain) compound had IC50 of 15 nM; the 2-(piperazinyl)benzimidazolyl variant had IC50 of 25 nM.

IT 307931-42-6P

RL: PUR (Purification or recovery); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of

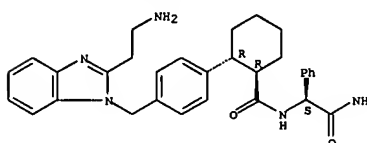
N-(phenylcyclohexylcarbonyl)phenyl

glycine amide for treatment of cardiovascular ischemia)

RN 307931-42-6 CAPLUS

CN Benzeneacetamide, α-([[(1R,2R)-2-[4-([2-(2-aminoethyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, dihydrochloride, (αS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



● 2 HCl

IT 307931-40-4P 307931-41-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(synthesis of heterocyclic derivs. of

N-(phenylcyclohexylcarbonyl)phenyl

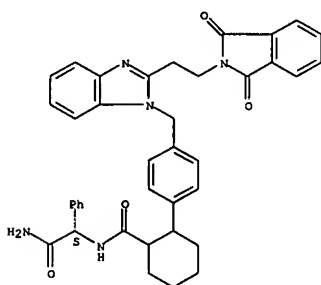
glycine amide for treatment of cardiovascular ischemia)

RN 307931-40-4 CAPLUS

CN Benzeneacetamide, α-([[(2-[4-([2-(1,3-dihydro-1,3-dioxo-2H-

isoindol-2-yl)ethyl]-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (αS)- (9CI) (CA INDEX NAME)

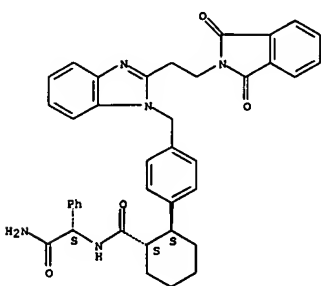
Absolute stereochemistry.



RN 307931-41-5 CAPLUS

CN Benzeneacetamide, α-([[(1S,2S)-2-[4-([2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)ethyl]-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (αS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



IT 307931-51-7P 307931-52-8P 307931-53-9P

307931-54-0P 307931-55-1P 307931-56-2P

307931-58-4P 307931-59-5P 307931-60-8P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of

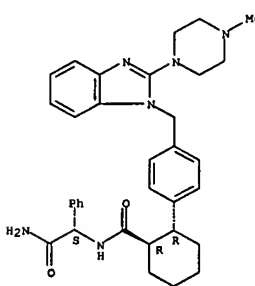
N-(phenylcyclohexylcarbonyl)phenyl

glycine amide for treatment of cardiovascular ischemia)

RN 307931-51-7 CAPLUS

CN Benzeneacetamide, α-([[(1R,2R)-2-[4-([2-(4-methyl-1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (αS)- (9CI) (CA INDEX NAME)

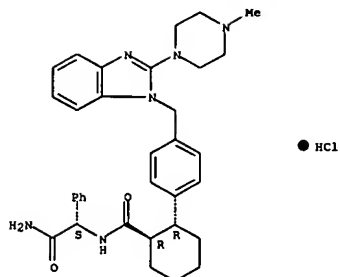
Absolute stereochemistry.



RN 307931-52-8 CAPLUS

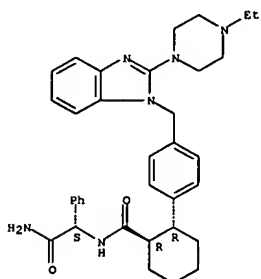
CN Benzeneacetamide, α-([[(1R,2R)-2-[4-([2-(4-methyl-1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, monohydrochloride, (αS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



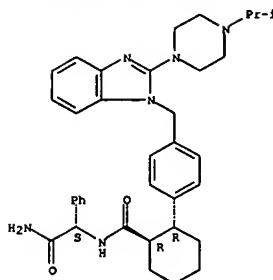
RN 307931-53-9 CAPLUS
 CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[[2-(4-ethyl-1-piperazinyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



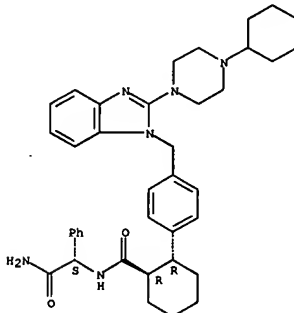
RN 307931-54-0 CAPLUS
 CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[[2-(4-(1-methylethyl)-1-piperazinyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



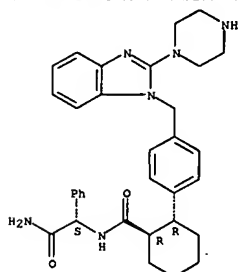
RN 307931-55-1 CAPLUS
 CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[[2-(4-cyclohexyl-1-piperazinyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



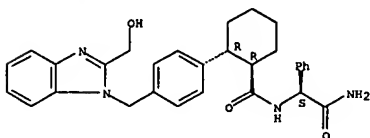
RN 307931-56-2 CAPLUS
 CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[[2-(1-piperazinyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



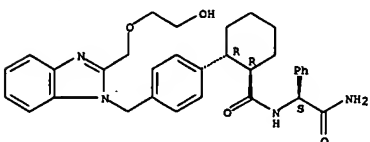
RN 307931-58-4 CAPLUS
 CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[[2-(2-hydroxymethyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

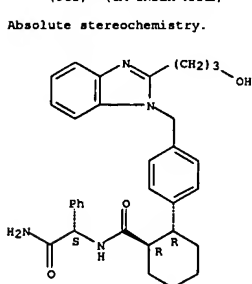


RN 307931-59-5 CAPLUS
 CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[[2-[(2-hydroxyethoxy)methyl]-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 307931-60-8 CAPLUS



Absolute stereochemistry.

L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 2000:841901 CAPLUS

DOCUMENT NUMBER: 133:362968

TITLE: Synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury

INVENTOR(S): Freund, Wolf-Dietrich; Lensky, Stephan; Muller, Stephan Nicholas; Paulsen, Holger; Keldenich, Jorg; Horvath, Ervin; Schuhmacher, Joachim

PATENT ASSIGNEE(S): Bayer A.-G., Germany

SOURCE: Ger. Offen., 30 pp. CODEN: GWKXBX

DOCUMENT TYPE: Patent

LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE

DE 19924818 A1 20001130 DE 1999-19924818 19990529

CA 2375186 AA 20001207 CA 2000-2375186 20000516

WO 2000073275 A1 20001207 WO 2000-EP4417 20000516

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, GR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA,

ZW RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

BR 2000011061 A 20020305 BR 2000-11061 20000516

EP 1185516 B1 20030502 EP 2000-925288 20000516

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO

JP 2003500475 T2 20030107 JP 2000-621341 20000516

EE 200100634 A 20030217 EE 2001-634 20000516

AT 238997 E 20030515 AT 2000-925288 20000516

AU 765752 B2 20030925 AU 2000-44055 20000516

PT 1185516 T 20030930 PT 2000-925288 20000516

ES 2197870 T3 20040116 ES 2000-925288 20000516

RU 2246490 C2 20050220 RU 2001-135715 20000516

ZA 2001009263 A 20021111 ZA 2001-9263 20011109

BG 106107 A 20020531 BG 2001-106107 20011113

NO 2001005810 A 20020125 NO 2001-5810 20011128

US 6716849 B1 20040406 US 2001-980243 20011129

HR 2001000955 A1 20030831 HR 2001-955 20011224

US 2004186106 A1 20040923 US 2004-805548 20040318

PRIORITY APPL. INFO.: DE 1999-19924818 A 19990529

WO 2000-EP4417 W 20000516

US 2001-980243 A3 20011129

OTHER SOURCE(S): MARPAT 133:362968

GI

L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

IT 307971-72-8P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307971-72-8 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

AB Title compds., e.g. I, were prepared for use in treating ischemic brain diseases in humans or animals. Thus I [X = N, X1 = O (II)] was prepared in six steps, starting from 2,3-diaminopyridine, glycolic acid, (1R,2R)-2-(4-bromomethylphenyl)cyclohexane-1-carboxylic acid tert-Bu ester (preparation given), and (S)-phenylglycine hydrochloride. Similarly prepared was I [X = C, X1 = N(Me) (III)]. In in vivo (binding of calf cortex adenosine transport protein) compds. II and III had Ki = 2 nM. In in vitro tests of rat brain reperfusion injury, II and III were effective at 0.001 mg/kg, reducing infarct volume 81-91% of control.

IT 307931-40-4

RL: RCT (Reactant); RACT (Reactant or reagent)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-40-4 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1,3-dioxo-2H-isoindol-2-yl)ethyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-58-4 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(hydroxymethyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 307931-59-5 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(2-hydroxyethoxy)methyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 307931-56-2P 307931-58-4P 307931-59-5P

307931-60-8P 307967-08-4P 307967-19-7P

307967-20-0P 307967-21-1P 307967-22-2P

307967-23-3P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(synthesis of heterocyclic derivs. of N-(phenylcyclohexylcarbonyl)phenylglycine amide for treatment of cerebral ischemia or injury)

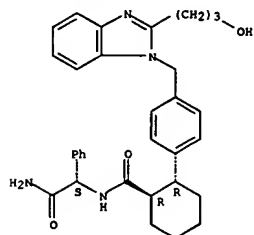
RN 307931-56-2 CAPLUS

CN Benzeneacetamide, α -[[(1R,2R)-2-[4-[(2-(1-piperazinyl)-1H-benzimidazol-1-yl)methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

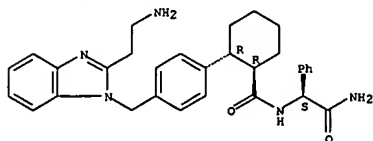
L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)
 RN 307931-60-8 CAPLUS
 CN Benzeneacetamide, α -[[[(1R,2R)-2-[4-[[2-(3-hydroxypropyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 307967-08-4 CAPLUS
 CN Benzeneacetamide, α -[[[(1R,2R)-2-[4-[[2-(2-aminoethyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

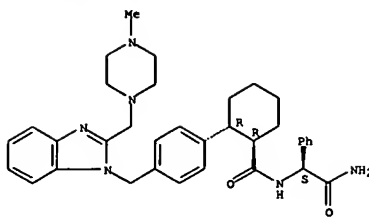
Absolute stereochemistry.



RN 307967-19-7 CAPLUS
 CN Benzeneacetamide, α -[[[(1R,2R)-2-[4-[[2-[(4-methyl-1-piperazinyl)methyl]-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

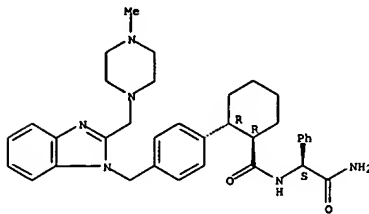
Absolute stereochemistry.

L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



RN 307967-20-0 CAPLUS
 CN Benzeneacetamide, α -[[[(1R,2R)-2-[4-[[2-[(4-methyl-1-piperazinyl)methyl]-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, monohydrochloride, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

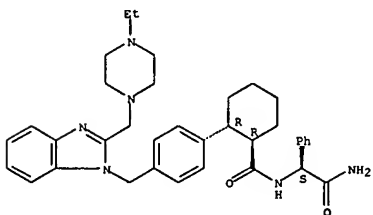


• HCl

RN 307967-21-1 CAPLUS
 CN Benzeneacetamide, α -[[[(1R,2R)-2-[4-[[2-[(4-ethyl-1-piperazinyl)methyl]-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

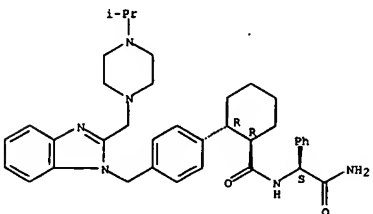
Absolute stereochemistry.

L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)



RN 307967-22-2 CAPLUS
 CN Benzeneacetamide, α -[[[(1R,2R)-2-[4-[[2-[(1-methylethyl)-1-piperazinyl]methyl]-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

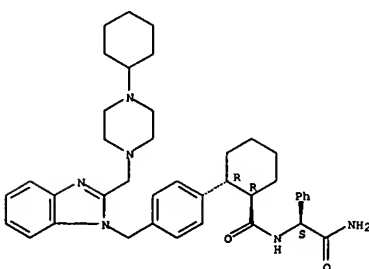
Absolute stereochemistry.



RN 307967-23-3 CAPLUS
 CN Benzeneacetamide, α -[[[(1R,2R)-2-[4-[[2-[(4-cyclohexyl-1-piperazinyl)methyl]-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]-, (α S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

L13 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN (Continued)

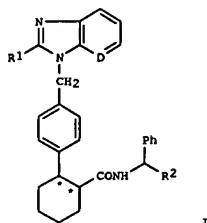


ACCESSION NUMBER: 1996:567333 CAPLUS
DOCUMENT NUMBER: 125:221843

TITLE: Preparation of benzylimidazole derivatives for the treatment of vascular restenosis
INVENTOR(S): Mueller-Gliemann, Matthias; Mueller, Ulrich; Beuck, Martin; Zaias, Siegfried; Gerdes, Christoph; Domdey-Bette, Anke; Gruetzmann, Rudi; Lohmer, Stefan; Wohlfeil, Stefan; et al.
PATENT ASSIGNEE(S): Bayer A.-G., Germany
SOURCE: Eur. Pat. Appl., 16 pp.
CODEN: EPKXDW
DOCUMENT TYPE: Patent
LANGUAGE: German
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

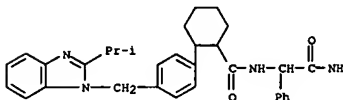
	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	EP 725064	A1	19960807	EP 1996-100760	19960119
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LI, LU, MC, NL, PT,				
SE	DE 19503160	A1	19960808	DE 1995-19503160	19950201
	TW 448176	B	20010801	TW 1996-85100684	19960122
	RO 117256	B1	20011228	RO 1996-152	19960126
	CA 2168317	AA	19960802	CA 1996-2168317	19960129
	JP 08253453	A2	19961001	JP 1996-33174	19960129
	IL 116931	A1	20000601	IL 1996-116931	19960129
	FI 9600425	A	19960802	FI 1996-425	19960130
	AU 9642240	A1	19960808	AU 1996-42240	19960130
	AU 710225	B2	19990916		
	BG 63044	B1	20010228	BG 1996-100326	19960130
	BG 103820	A	20010928	BG 1999-103820	19960130
	NO 9600414	A	19960802	NO 1996-414	19960131
	NO 309841	B1	20010409		
	ZA 9600725	A	19960820	ZA 1996-725	19960131
	RU 2158261	C2	20001027	RU 1996-101800	19960131
	CN 1137380	A	19961211	CN 1996-102574	19960201
	US 5935963	A	19990810	US 1997-960075	19971024
PRIORITY APPL. INFO.:				DE 1995-19503160	A 19950201
				US 1996-588477	B1 19960118

OTHER SOURCE(S): MARPAT 125:221843
GI

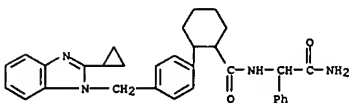


I

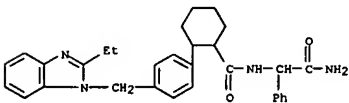
AB The title compds. [I: D = CH, N; R1 = Ph, cycloalkyl, (un)branched alkyl; R2 = (un)branched alkoxy carbonyl, CH2OH, CONH2], useful for the treatment of vascular restenosis, are prepared. Thus, I (D = N, R1 = CHMe2, R2 = CONH2; * * cyclohexyl ring bonding is trans) was prepared and demonstrated a IC50 of 0.01 nM for the inhibition of rat aorta smooth muscle proliferation.
IT 181130-37-OP 181130-38-1P 181130-41-6P
181130-42-7P
RL: EAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of benzylimidazole derivs. for the treatment of vascular restenosis)
RN 181130-37-0 CAPLUS
CN Benzeneacetamide, α-[[[2-[4-[[2-(1-methylethyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]- (9CI) (CA INDEX NAME)



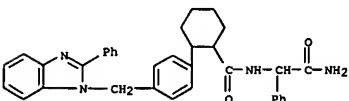
RN 181130-38-1 CAPLUS
CN Benzeneacetamide, α-[[[2-[4-[[2-(cyclopropyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]- (9CI) (CA INDEX NAME)



RN 181130-41-6 CAPLUS
CN Benzeneacetamide, α-[[[2-[4-[[2-(ethyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]- (9CI) (CA INDEX NAME)



RN 181130-42-7 CAPLUS
CN Benzeneacetamide, α-[[[2-[4-[[2-(phenyl)-1H-benzimidazol-1-yl]methyl]phenyl]cyclohexyl]carbonyl]amino]- (9CI) (CA INDEX NAME)



=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

15.79

522.58

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

-2.25

-2.25

STN INTERNATIONAL LOGOFF AT 10:06:55 ON 17 AUG 2006